CLAIMS

What is claimed is:

5 1. A compound having the formula

$$\begin{array}{c} B \\ Y \\ \stackrel{\stackrel{\longrightarrow}{=}}{=} \\ R_3 \end{array} \qquad \begin{array}{c} N \\ N \\ N \\ N \end{array} \qquad \begin{array}{c} R_2 \\ N \\ N \\ N \end{array} \qquad \begin{array}{c} R_2 \\ N \\ N \\ N \end{array} \qquad \begin{array}{c} (I) \\ N \\ N \\ N \end{array}$$

wherein:

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- (a) R₁ is C₁₋₈ alkyl, C₃₋₇ cycloalkyl, or C₄₋₁₀ (alkylcycloalkyl), which are all optionally substituted from one to three times with halo, cyano, nitro, C₁₋₆ alkoxy, amido, amino or phenyl, or R₁ is C₆ or C₁₀ aryl which is optionally substituted from one to three times with halo, cyano, nitro, C₁₋₆ alkyl, C₁₋₆ alkoxy, amido, amino or phenyl;
- (b) m is 1 or 2;
 - (c) n is 1 or 2;
 - (d) R₂ is C₁₋₆ alkyl, C₂₋₆ alkenyl or C₃₋₇ cycloalkyl, each optionally substituted from one to three times with halogen, or R₂ is H;
 - (e) R_3 is C_{1-8} alkyl optionally substituted with phenyl, C_{3-12} alkenyl, C_{3-7} cycloalkyl, or C_{4-10} (alkylcycloalkyl), wherein the cycloalkyl or

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(h)

alkylcycloalkyl are optionally substituted with hydroxy, C_{1-6} alkyl, C_{2-6} alkenyl; or C_{1-6} alkoxy or R_3 together with the carbon atom to which it is attached forms a C_{3-7} cycloalkyl group optionally substituted with C_{2-6} alkenyl;

(f) Y is H, phenyl substituted with nitro, pyridyl substituted with nitro, or C_{1-6} alkyl wherein said alkyl is optionally substituted with cyano, OH or C_{3-7} cycloalkyl;

 R_4 is (i) C_{1-10} alkyl optionally substituted with

- 10 (g) B is H, C_{1-6} alkyl, $R_4-(C=0)-$, $R_4O(C=0)-$, $R_4-N(R_5)-C(=0)-$, $R_4-N(R_5)-C(=S)-$, R_4SO_2- , or $R_4-N(R_5)-SO_2-$;
 - phenyl, carboxyl, C1-6 alkanovl, 1-3 halogen. hydroxy, -OC(0)C1-6 alkyl, C1-6 alkoxy, amino optionally mono-or-di substituted with C1-6 alkyl, amido, or (lower alkyl) amido; or -O-phenyl optionally substituted with halogen or C1-6 alkoxy; (ii) C_{3-7} cycloalkyl, C_{3-7} cycloalkoxy, or C4-10 alkylcyclo-alklyl, all optionally substituted with hydroxy, carboxyl, (C1-6 alkoxy) carbonyl, amino optionally mono- or disubstituted with C_{1-6} alkyl, amido, or (lower alkyl) amido; (iii) amino optionally mono-or-di-substituted with C1-6 alkyl; amido; or (lower alkyl)amido; (iv) C6 or C10 aryl or C_{7-16} aralkyl, all optionally substituted with C1-6 alkyl, halogen, nitro, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-disubstituted with C_{1-6} alkyl; or (v) Het or (lower alkyl)-Het, both optionally substituted with C1-6

alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C1-6

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alkyl; (vi) bicyclo(1.1.1)pentane; (vii) $-C(0)OC_{1-6}$ alkyl, C_{26} alkenyl, C_{26} alkynyl; and

- (i) R_5 is H or C_{1-6} alkyl, said C_{1-6} alkyl optionally substituted with 1-3 halogens;
- or a pharmaceutically acceptable salt, solvate or prodrug thereof.
 - 2. A compound of Claim 1 wherein m is 2.
- 10 3. A compound of Claim 1 wherein n is 1.
 - 4. A compound of Claim 1 wherein R1 is cyclopropyl.
 - 5. A compound of Claim 1 wherein R₁ is cyclobuty1.
 - 6. A compound of Claim 1 wherein R_1 is optionally substituted phenyl.
- 7. A compound of Claim 1 wherein R_2 is ethyl or 20 vinyl.
 - 8. A compound of Claim 1 wherein R_3 is C_{1-6} alkyl.
- 9. A compound of Claim 1 wherein m is 2, n is 1 and $R_2 \mbox{ is ethy1.}$
 - 10. A compound of Claim 9 wherein R_1 is cyclopropy1.
 - 11. A compound of Claim 9 wherein R_1 is cyclobutyl.
 - 12. A compound of Claim 9 wherein R_1 is optionally substituted phenyl.

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- 13. A compound of Claim 1 wherein m is 2, n is 1 and R_2 is vinyl.
- 14. A compound of Claim 13 wherein R_1 is cyclopropyl.
- 15. A compound of Claim 13 wherein R1 is cyclobutyl.
- 16. A compound of Claim 13 wherein R_1 is optionally substituted phenyl.

17. A compound having the formula

$$\begin{array}{c} \text{B} \\ \text{Y} \\ \stackrel{\stackrel{\circ}{=}}{\overset{\circ}{\mathbb{R}}_{3}} \end{array}$$

wherein:

- (a) R₁₁ is C₁₋₈ alkyl, C₃₋₇ cycloalkyl, or
 C₄₋₁₀(alkylcyclo-alkyl), naphthyl, or phenyl wherein said phenyl is optionally substituted from one to three times with halo, cyano, nitro, C₁₋₆ alkyl, C₁₋₆ alkoxy, amido, or phenyl;
 - (b) R₁₂ is C₁₋₆ alkyl, C₂₋₆ alkenyl or H;
- 20 (c) R₃ is C₁₋₈ alkyl, C₃₋₁₂ alkenyl, C₃₋₇ cycloalkyl, or C₄₋₁₀ (alkylcycloalkyl), wherein the cycloalkyl or alkylcycloalkyl are optionally substituted with hydroxy, C₁₋₆ alkyl, C₁₋₆ alkenyl, or C₁₋₆ alkoxy;

- (d) Y is H or C₁₋₆ alkyl wherein said alkyl is optionally substituted with cyano or C₃₋₇ cycloalkyl;
- (e) B is H, R_4 -(C=O)-, R_4 O(C=O)-, R_4 -N(R_5)-C(=O)-, R_4 -N(R_5)-C(=S)-, R_4 SO₂-, or R_4 -N(R_5)-SO₂-;
- (f) R_4 is (i) C_{1-10} alkyl optionally substituted with carboxyl, C_{1-6} alkanoyl, hydroxy, C_{1-6} alkoxy, amino optionally mono-or-di substituted with C_{1-6} alkyl, amido, or (lower alkyl) amido; (ii) C_{3-7}
- cycloalkyl, C₃₋₇ cycloalkoxy, or C₄₋₁₀ alkylcycloalklyl, all optionally substituted with hydroxy, carboxyl, (C₁₋₆ alkoxy)carbonyl, amino optionally mono- or disubstituted with C₁₋₆ alkyl, amido, or (lower alkyl) amido; (iii) amino optionally monoor-di-substituted with C₁₋₆ alkyl, amido; or (lower alkyl)amido; (iv) C₆ or C₁₀ aryl or C₇₋₁₆ aralkyl, all optionally substituted with C₁₋₆ alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C₁₋₆ alkyl;
- or (v) Het or (lower alkyl)-Het, both optionally substituted with C₁₋₆ alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-disubstituted with C₁₋₆ alkyl; and
 - (g) R_5 is H or C_{1-6} alkyl,
- 25 or a pharmaceutically acceptable salt, solvate or prodrug thereof.
 - 18. A compound of Claim 17 wherein R₁₁ is selected from cyclopropyl, cyclobutyl or optionally substituted phenyl.
 - 19. A compound having the formula

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wherein:

- (a) R₁₁ is C₁₋₈ alkyl, C₃₋₇ cycloalkyl, or C₄₋₁₀ (alkylcyclo-alkyl), naphthyl, or phenyl wherein said phenyl is optionally substituted from one to three times with halo, cyano, nitro, C₁₋₆ alkyl, C₁₋₆ alkoxy, amido, or phenyl;
- (b) R₃ is C₁₋₈ alkyl, C₃₋₁₂ alkenyl, C₃₋₇ cycloalkyl, or C₄₋₁₀ (alkylcycloalkyl), wherein the cycloalkyl or alkylcycloalkyl are optionally substituted with hydroxy, C₁₋₆ alkyl, C₁₋₆ alkenyl, or C₁₋₆ alkoxy;
- Y is H or C₁₋₆ alkyl wherein said alkyl is optionally substituted with cyano or C₃₋₇ cycloalkyl;
- 15 (d) B is H, R_4 -(C=O)-, R_4 O(C=O)-, R_4 -N(R_5)-C(=O)-, R_4 -N(R_5)-C(=S)-, R_4 SO₂-, or R_4 -N(R_5)-SO₂-;
- (e) R₄ is (i) C₁₋₁₀ alkyl optionally substituted with carboxyl, C₁₋₆ alkanoyl, hydroxy, C₁₋₆ alkoxy, amino optionally mono-or-di substituted with C₁₋₆ alkyl,
 amido, or (lower alkyl) amido; (ii) C₃₋₇ cycloalkyl, C₃₋₇ cycloalkoxy, or C₄₋₁₀ alkylcycloalkyl, all optionally substituted with hydroxy, carboxyl, (C₁₋₆ alkoxy)carbonyl, amino optionally mono- or disubstituted with C₁₋₆ alkyl, amido, or

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(lower alkyl) amido; (iii) amino optionally mono-or-di-substituted with C_{1-6} alkyl; amido; or (lower alkyl) amido; (iv) C_6 or C_{10} aryl or C_{7-16} aralkyl, all optionally substituted with C_{1-6} alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C_{1-6} alkyl; or (v) Het or (lower alkyl)-Het, both optionally substituted with C_{1-6} alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C_{1-6} alkyl; and

- (f) R_{5} is H or $C_{1\text{-}6}$ alkyl; or a pharmaceutically acceptable salt, solvate or prodrug thereof.
- 15 20. A compound of Claim 19 wherein R_{11} is selected from cyclopropyl, cyclobutyl or optionally substituted phenyl.
 - 21. A compound having the formula

wherein:

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 (a) R₁₁ is C₁₋₈ alkyl, C₃₋₇ cycloalkyl, or C₄₋₁₀ (alkylcyclo-alkyl), naphthyl, or phenyl

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wherein said phenyl is optionally substituted from one to three times with halo, cyano, nitro, C_{1-6} alkyl, C_{1-6} alkoxy, amido, or phenyl;

- (b) R₃ is C₁₋₈ alkyl, C₃₋₁₂ alkenyl, C₃₋₇ cycloalkyl, or C₄₋₁₀ (alkylcycloalkyl), wherein the cycloalkyl or alkylcycloalkyl are optionally substituted with hydroxy, C₁₋₆ alkyl, C₁₋₆ alkenyl, or C₁₋₆ alkoxy;
 - (c) Y is H or C₁₋₆ alkyl wherein said alkyl is optionally substituted with cyano or C₃₋₇ cycloalkyl;
 - (d) B is H, R_4 -(C=O)-, R_4 O(C=O)-, R_4 -N(R_5)-C(=O)-, R_4 -N(R_5)-C(=S)-, R_4 SO₂-, or R_4 -N(R_5)-SO₂-;
- (e) R₄ is (i) C₁₋₁₀ alkyl optionally substituted with carboxyl, C₁₋₆ alkanoyl, hydroxy, C₁₋₆ alkoxy, amino optionally mono-or-di substituted with C₁₋₆ alkyl, amido, or (lower alkyl) amido; (ii) C₃₋₇ cycloalkyl, C₃₋₇ cycloalkoxy, or C₄₋₁₀ alkylcycloalkyl, all optionally substituted with hydroxy, carboxyl, (C₁₋₆ alkoxy)carbonyl, amino optionally mono- or disubstituted with C₁₋₆ alkyl, amido, or (lower alkyl) amido; (iii) amino optionally mono- or-di-substituted with C₁₋₆ alkyl; amido; or (lower alkyl)amido; (iv) C₆ or C₁₀ aryl or C₇₋₁₆ aralkyl, all optionally substituted with C₁₋₆ alkyl, hydroxy, amido, (lower alkyl) amido, or amino
- 25 hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C₁₋₆ alkyl; or (v) Het or (lower alkyl)-Het, both optionally substituted with C₁₋₆ alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C₁₋₆ alkyl; and
 - (f) R_5 is H or C_{1-6} alkyl; or a pharmaceutically acceptable salt, solvate or prodrug thereof.

- 22. A compound of Claim 21 wherein R₁₁ is selected from cyclopropyl, cyclobutyl or optionally substituted phenyl.
- 23. A compound having the formula

$$\begin{array}{c} \text{B} \\ \text{Y} \\ \stackrel{\stackrel{\bullet}{=}}{\overset{\bullet}{\mathbb{R}}_{3}} \end{array}$$

wherein:

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- (a) R_{11} is C_{1-8} alkyl, C_{3-7} cycloalkyl, or 10 C_{4-10} (alkylcycloalkyl), naphthyl, or phenyl wherein said phenyl is optionally substituted from one to three times with halo, cyano, nitro, C_{1-6} alkyl, C_{1-6} alkoxy, amido, or phenyl;
 - (b) R₃ is C₁₋₈ alkyl, C₃₋₁₂ alkenyl, C₃₋₇ cycloalkyl, or C₄₋₁₀ (alkylcycloalkyl), wherein the cycloalkyl or alkylcycloalkyl are optionally substituted with hydroxy, C₁₋₆ alkyl, C₁₋₆ alkenyl, or C₁₋₆ alkoxy;
 - (c) Y is H or C₁₋₆ alkyl wherein said alkyl is optionally substituted with cyano or C₃₋₇ cycloalkyl;
 - (d) B is H, R_4 -(C=O)-, R_4 O(C=O)-, R_4 -N(R_5)-C(=O)-, R_4 -N(R_5)-C(=S)-, R_4 SO₂-, or R_4 -N(R_5)-SO₂-;
 - (e) R_4 is (i) $C_{1\text{--}10}$ alkyl optionally substituted with

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carboxyl, C1-6 alkanoyl, hydroxy, C1-6 alkoxy, amino optionally mono-or-di substituted with C1-6 alkyl, amido, or (lower alkyl) amido; (ii) C3-7 cycloalkyl, C3-7 cycloalkoxy, or C4-10 alkylcycloalklyl, all optionally substituted with hydroxy, carboxyl, (C1-6 alkoxy) carbonyl, amino optionally mono- or disubstituted with C1-6 alkyl, amido, or (lower alkyl) amido; (iii) amino optionally monoor-di-substituted with C1-6 alkyl; amido; or (lower alkyl) amido; (iv) C6 or C10 aryl or C7-16 aralkyl, all optionally substituted with C1-6 alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C1-6 alkyl; or (v) Het or (lower alkyl)-Het, both optionally substituted with C1-6 alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-disubstituted with C1-6 alkyl;

- (f) R₅ is H or C₁₋₆ alkyl;
- (g) n is 1 or 2; and
- 20 (h) p is 1, 2, 3, 4 or 5, or a pharmaceutically acceptable salt, solvate or prodrug thereof.
- 24. A compound of Claim 23 wherein R₁₁ is selected from cyclopropyl, cyclobutyl or optionally substituted phenyl.
 - 25. A compound of having the formula

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wherein:

- (a) R₃₁ is C₁₋₈ alkyl, C₃₋₇ cycloalkyl, or C₄₋₁₀ (alkylcycloalkyl), all optionally substituted with hydroxy, halo, C₁₋₆ alkoxy, C₁₋₆ thioalkyl, amido, amino, (C₁₋₆ alkyl)amido, C6 or C₁₀ aryl, C₇₋₁₆ aralkyl, Het, or (C₁₋₆ alkyl)-Het, said aryl, arylalkyl or Het being optionally substituted with halo, alkyl or lower alkyl Het;
- 10 (b) n is 1 or 2;
 - (c) R₃₂ is H, C₁₋₆ alkyl, C₁₋₃ alkoxy, C₃₋₇ cycloalkyl, C₂₋₆ alkenyl, or C₂₋₆ alkynyl, all optionally substituted with halogen;
 - (d) R₁₃ is C₁₋₈ alkyl, C₃₋₁₂ alkenyl, C₃-C₇ cycloalkyl, C₄₋₁₃ cycloalkenyl, or C₄-C₁₀ (alkylcycloalkyl), all optionally substituted with hydroxy, C₁-C₆ alkoxy, C₁-C₆ thioalkyl, amino, amido, (loweralkyl) amido, C₆ or C₁₀ aryl, or C₇-C₁₆ aralkyl;
 - (e) Y2 is H or C1-C6 alkyl;
- 20 (f) B_2 is H, R_{14} -(C=O)-; R_{14} O(C=O)-, R_{14} -N(R_{15})-C(=O)-; R_{14} -N(R_{15})-C(=S)-; R_{14} SO₂-, or R_{14} -N(R_{15})-SO₂-;
 - (g) R_{14} is (i) C_{1-10} alkyl optionally substituted with carboxyl, C_{1-6} alkanoyl, hydroxy, C_{1-6} alkoxy, amino

optionally mono-or-di substituted with C1-6 alkyl, amido, or (lower alkyl) amido; (ii) C3-7 cycloalkyl, C3-7 cycloalkoxy, or C4-10 alkylcycloalkyl, all optionally substituted with hydroxy, carboxyl, (C1-6 alkoxy) carbonyl, amino optionally mono- or disubstituted with C1-6 alkyl, amido, or (lower alkyl) amido; (iii) amino optionally mono-or-di-substituted with C1-6 alkyl; amido; or (lower alkyl)amido; (iv) C6 or C10 aryl or C7-16 aralkyl, all optionally substituted with C1.6 alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C1-6 alkyl; or (v) Het or (lower alkyl)-Het, both optionally substituted with C1-6 alkyl, hydroxy, amido, (lower alkyl) amido, or amino optionally mono-or-di-substituted with C1-6 alkyl; and

- (h) R_{15} is H or C_{1-6} alkyl.
- A salt, solvate or prodrugs of a compound of
 Claim 25.
 - 27. A compound of Claim 25 wherein $R_{31} \text{ is } C_{3-6} \text{ cycloalkyl}, \ C_{4^-10} \text{ alkylcycloalkyl}, \\ C_{1-8} \text{ alkyl } CF_3 \text{ or } CCl_3.$

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- 28. A compound of Claim 25 wherein B_2 is an acyl derivative of formula R_{14} -O-(C=O)- or a carboxyl of formula R_{14} -O-(C=O)-.
- 30 29. A compound of claim 25 wherein R_2 is H, C_{1-3} alkyl, C_{3-5} cycloalkyl, or C_{2-4} alkenyl, all optionally substituted with halo.

30. A compound of claim 25 wherein R_{31} is C_{1-8} alkyl, C_{3-7} cycloalkyl, or C_{4-10} alkylcycloalkyl, all optionally substituted with hydroxy, C_{1-6} alkoxy, C_{1-6} thioalkyl, acetamido or C_6 or C_{10} aryl.

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31. A compound of claim 25 wherein
B is (CH₃)₃-O-CO-;
Y is H; n is 1;
R₃₁ is methyl, cyclopropyl or -CF₃;
R₃₂ is ethyl or vinyl; and
R₁₃ is t-butyl, i-propyl, s-butyl, i-butyl or

cyclohexylmethyl.

- 32. A pharmaceutical composition, comprising
 - (a) a compound of Claim 1-31, or a pharmaceutically acceptable salt, solvate or prodrug thereof; and
 - (b) a pharmaceutically acceptable carrier.
- 20 33. A method of inhibiting HCV NS3 protease which comprises administering to a mammal in need of such treatment a therapeutically effective amount a compound of Claim 1-31, or a pharmaceutically acceptable salt, solvate or prodrug thereof.

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34. A method of for treating an HCV infection, in a mammal in need thereof, comprising the administration to said mammal of a therapeutically effective amount a compound of Claim 1-31, or a pharmaceutically acceptable salt, solvate or prodrug thereof.